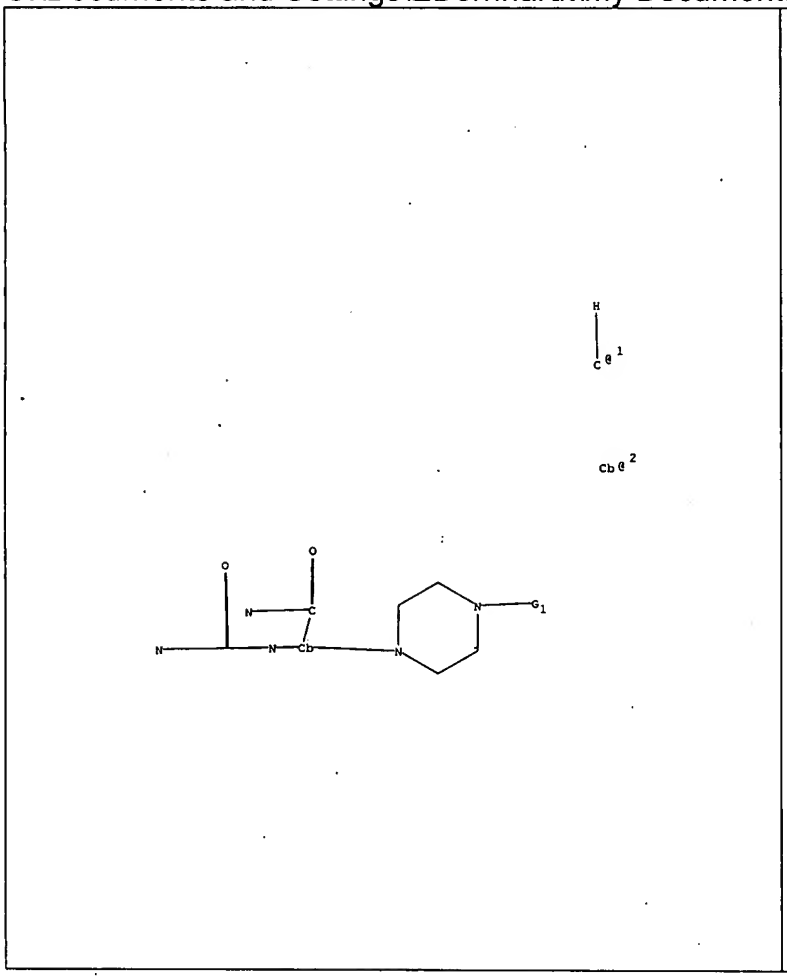
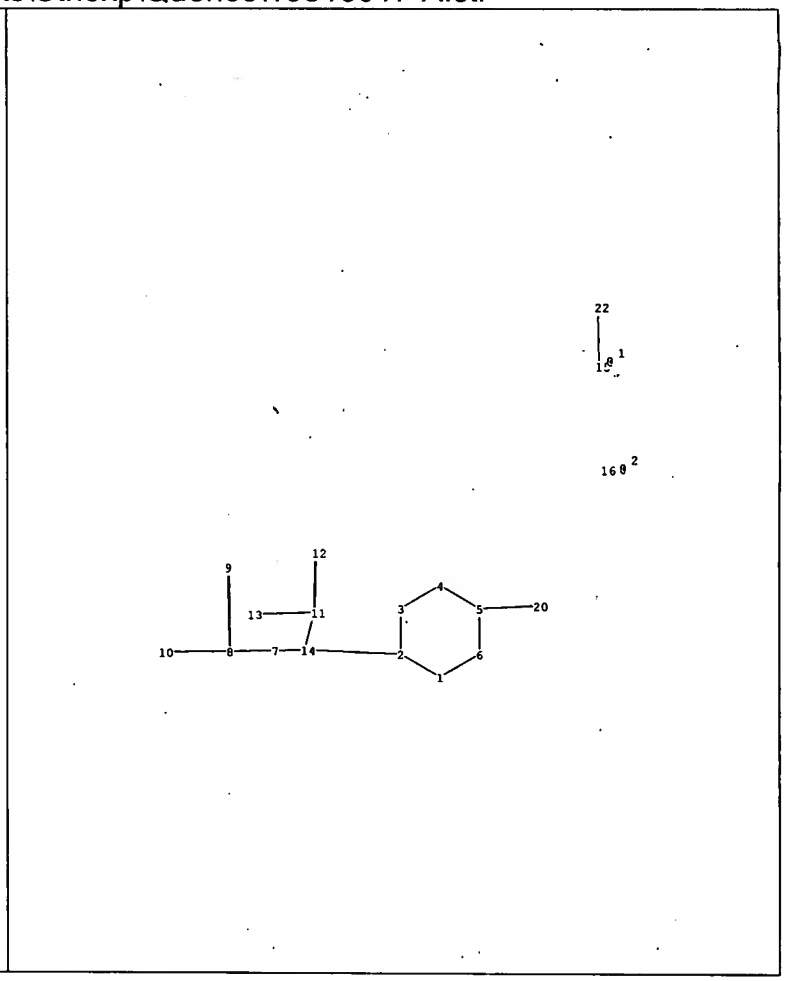
EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L4	3318	((544/357) or (544/377) or (544/381) or (544/396) or (544/400) or (540/575)).CCLS.	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:35
L5	48	1-piperazinyl near25 benzamide	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:36
L6	13	l4 and l5	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:37
L7	35	l5 not l6	US-PGPUB; USPAT	OR	OFF	2007/04/28 18:37

C:\Documents and Settings\EBernhardt\My Documents\Stnexp\Queries\10815017-A.str





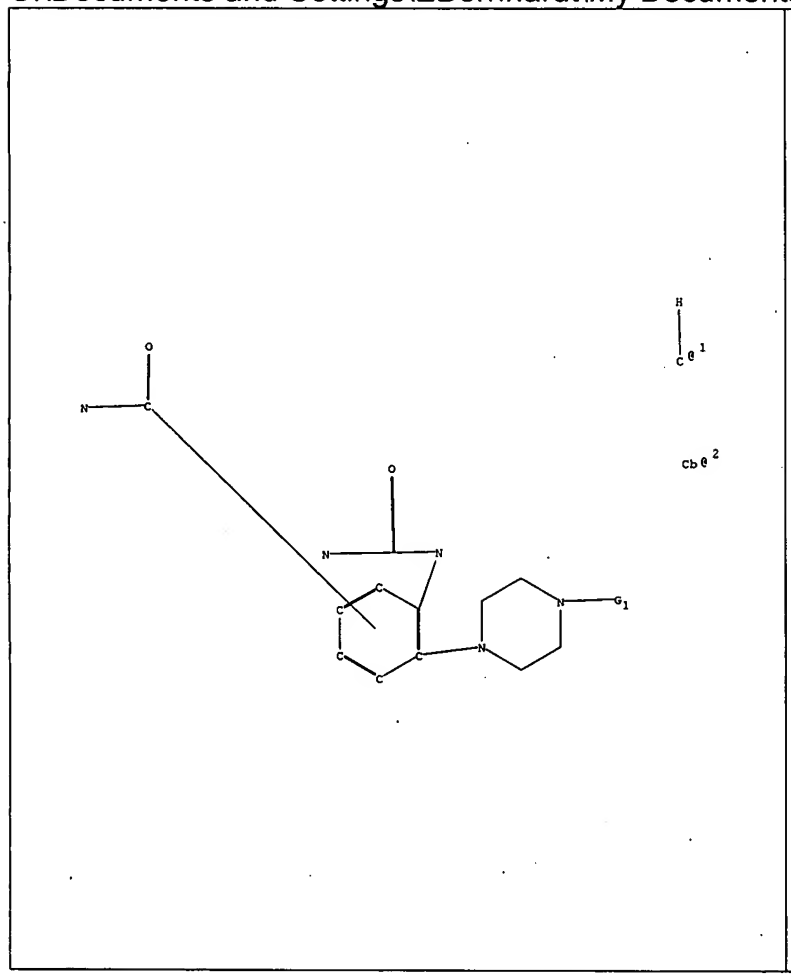
chain nodes :
7 8 9 10 11 12 14 15 16 20 22
ring nodes :
1 2 3 4 5 6
ring/chain nodes :
13
chain bonds :
2-14 5-20 7-8 7-14 8-9 8-10 11-12 11-13 11-14 15-22
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-20 7-8 8-9 8-10 11-12 11-13
exact bonds :
2-14 7-14 11-14 15-22
isolated ring systems :
containing 1 :

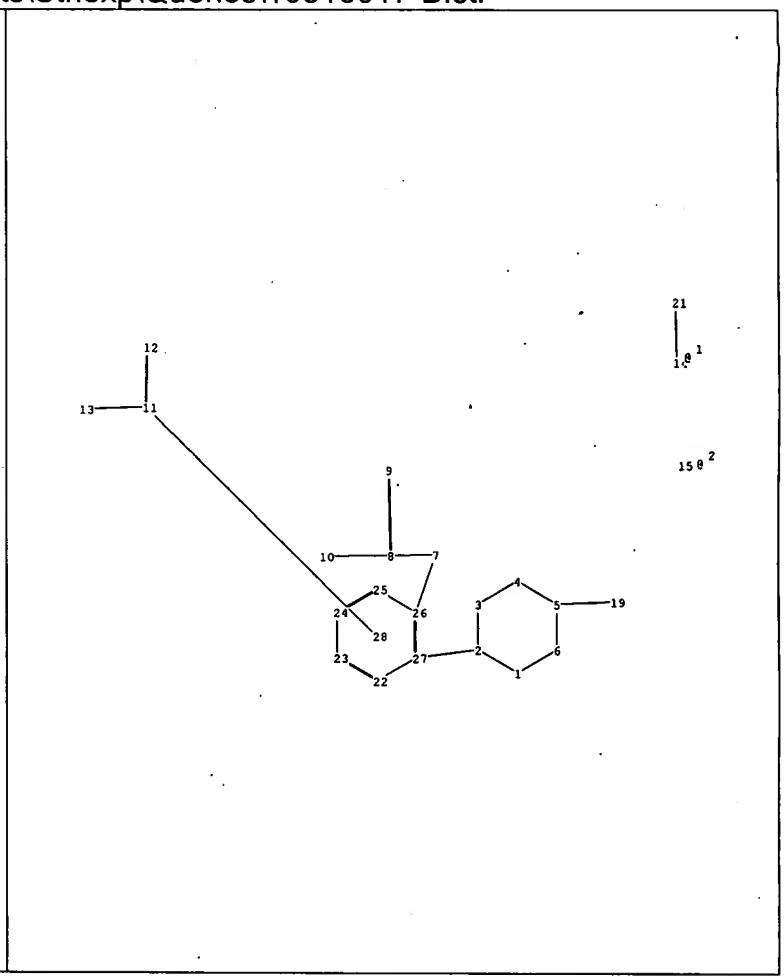
G1:[*1],[*2]

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS
12:CLASS 13:CLASS 14:Atom 15:CLASS 16:Atom 20:CLASS 22:CLASS
Generic attributes :

14:
Saturation : Unsaturated
Number of Carbon Atoms : less than 7
Type of Ring System : Monocyclic

C:\Documents and Settings\EBernhardt\My Documents\Stnexp\Queries\10815017-B.str





chain nodes :
7 8 9 10 11 12 14 15 19 21
ring nodes :
1 2 3 4 5 6 22 23 24 25 26 27
ring/chain nodes :
13
chain bonds :
2-27 5-19 7-8 7-26 8-9 8-10 11-12 11-13 14-21
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 22-23 22-27 23-24 24-25 25-26 26-27
exact/norm bonds :
1-2 1-6 2-3 2-27 3-4 4-5 5-6 5-19 7-8 7-26 8-9 8-10 11-12 11-13
exact bonds :
14-21
normalized bonds :
22-23 22-27 23-24 24-25 25-26 26-27
isolated ring systems :
containing 1 :

G1:[*1],[*2]

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS
12:CLASS

13:CLASS 14:CLASS 15:Atom 19:CLASS 21:CLASS 22:CLASS 23:Atom 24:Atom 25:Atom 26:CLASS 27:Atom 28:Atom

```
=> s l1
SAMPLE SEARCH INITIATED 16:08:07 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     1916 TO ITERATE

100.0% PROCESSED     1916 ITERATIONS                                 34 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        35695 TO    40945
PROJECTED ANSWERS:             331 TO     1029

L2            34 SEA SSS SAM L1

=> d l2 1 5 10 30

L2   ANSWER 1 OF 34  REGISTRY  COPYRIGHT 2007 ACS on STN
RN   861982-63-0  REGISTRY
ED   Entered STN:  29 Aug 2005
CN   Benzamide, N-(3-methoxypropyl)-2-[4-(phenylmethyl)-1-piperazinyl]-5-[[[[2-
     (trifluoromethyl)phenyl]amino]carbonyl]amino]- (9CI)  (CA INDEX NAME)
MF   C30 H34 F3 N5 O3
SR   Chemical Library
       Supplier: ComGenex International Inc.
LC   STN Files:   CHEMCATS
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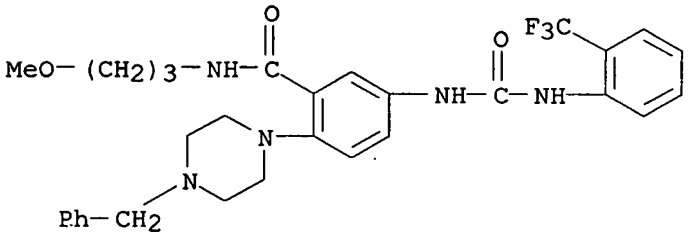


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**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**


L2   ANSWER 5 OF 34  REGISTRY  COPYRIGHT 2007 ACS on STN
RN   860078-73-5  REGISTRY
ED   Entered STN:  12 Aug 2005
CN   Pyrrolidine, 1-[5-[[[(3,5-dimethoxyphenyl)amino]carbonyl]amino]-2-[4-(2-
     methoxyphenyl)-1-piperazinyl]benzoyl]- (9CI)  (CA INDEX NAME)
MF   C31 H37 N5 O5
SR   Chemical Library
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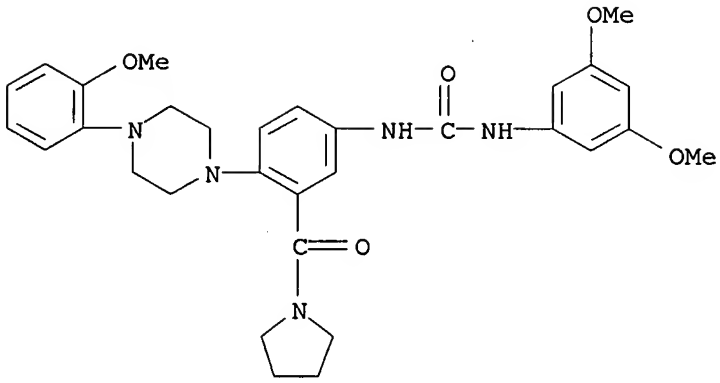
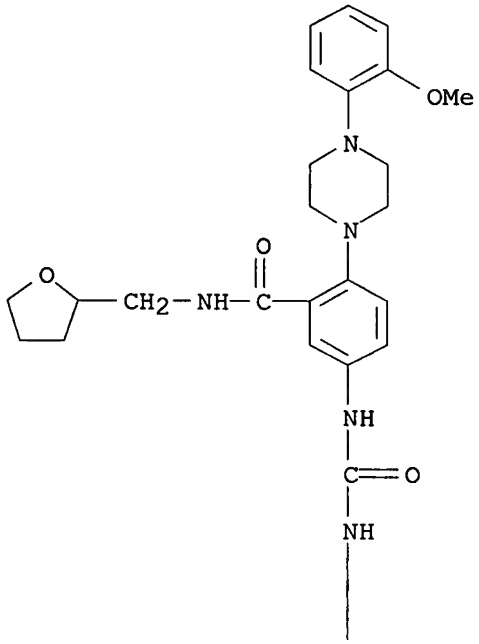


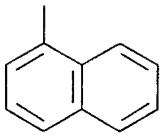
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 ANSWER 10 OF 34 REGISTRY COPYRIGHT 2007 ACS on STN
RN 857668-00-9 REGISTRY
ED Entered STN: 29 Jul 2005
CN Benzamide, 2-[4-(2-methoxyphenyl)-1-piperazinyl]-5-[[(1-naphthalenylamino)carbonyl]amino]-N-[(tetrahydro-2-furanyl)methyl]- (9CI) (CA INDEX NAME)
MF C34 H37 N5 O4
SR Chemical Library
Supplier: ComGenex International Inc.
LC STN Files: CHEMCATS

PAGE 1-A



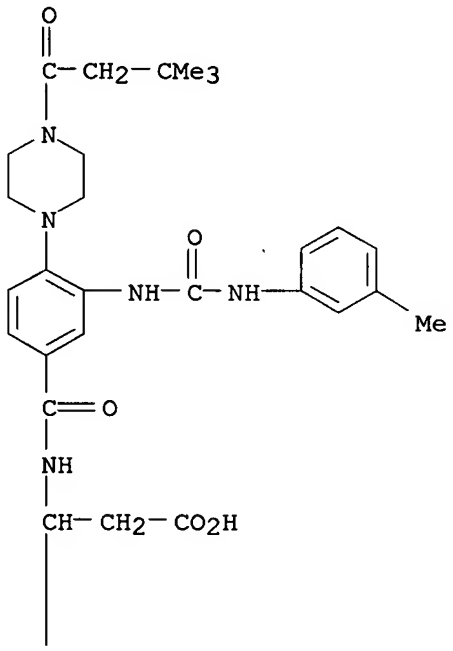
PAGE 2-A

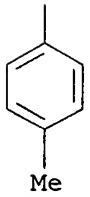


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 ANSWER 30 OF 34 REGISTRY COPYRIGHT 2007 ACS on STN
RN 439274-60-9 REGISTRY
ED Entered STN: 18 Jul 2002
CN Benzenepropanoic acid, β-[[4-[4-(3,3-dimethyl-1-oxobutyl)-1-piperazinyl]-3-[[[(3-methylphenyl)amino]carbonyl]amino]benzoyl]amino]-4-methyl- (9CI) (CA INDEX NAME)
MF C35 H43 N5 O5
SR Chemical Library
Supplier: Ambinter

PAGE 1-A





PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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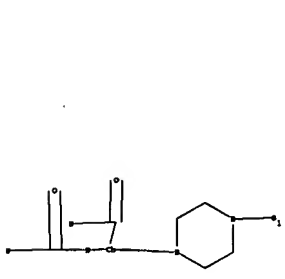
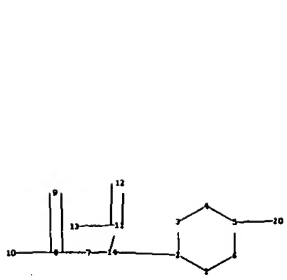
Uploading C:\Documents and Settings\EBernhardt\My Documents\Stnexp\Queries\10815017-A.str





chain nodes :
7 8 9 10 11 12 14 15 16 20 22
ring nodes :
1 2 3 4 5 6
ring/chain nodes :
13

```
chain bonds :
2-14  5-20  7-8  7-14  8-9  8-10  11-12  11-13  11-14  15-22
ring bonds :
1-2  1-6  2-3  3-4  4-5  5-6
exact/norm bonds :
1-2  1-6  2-3  3-4  4-5  5-6  5-20  7-8  8-9  8-10  11-12  11-13
exact bonds :
2-14  7-14  11-14  15-22
isolated ring systems :
containing 1 :


G1:[*1],[*2]

Match level :
1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:CLASS  8:CLASS  9:CLASS  10:CLASS
 11:CLASS  12:CLASS  13:CLASS  14:Atom  15:CLASS  16:Atom  20:CLASS  22:CLASS
Generic attributes :
14:
Saturation              : Unsaturated
Number of Carbon Atoms : less than 7
Type of Ring System     : Monocyclic



L3      STRUCTURE UPLOADED

=> s l3
SAMPLE SEARCH INITIATED 16:10:14 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     1916 TO ITERATE

100.0% PROCESSED     1916 ITERATIONS                                  27 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        35695 TO    40945
PROJECTED ANSWERS:             229 TO      851

L4           27 SEA SSS SAM L3

=> d l4 1 5 10 27

L4   ANSWER 1 OF 27  REGISTRY  COPYRIGHT 2007 ACS on STN
RN   861982-63-0  REGISTRY
ED   Entered STN:  29 Aug 2005
CN   Benzamide, N-(3-methoxypropyl)-2-[4-(phenylmethyl)-1-piperazinyl]-5-[[[[2-
     (trifluoromethyl)phenyl]amino]carbonyl]amino]- (9CI)  (CA INDEX NAME)
MF   C30 H34 F3 N5 O3
SR   Chemical Library
       Supplier: ComGenex International Inc.
LC   STN Files:   CHEMCATS
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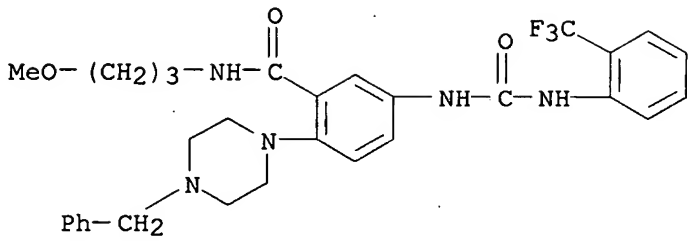
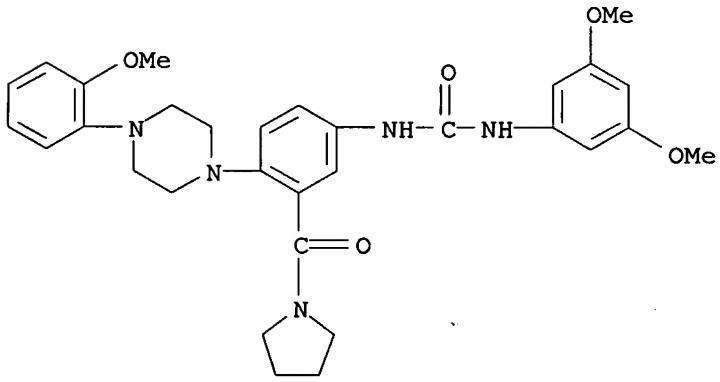


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

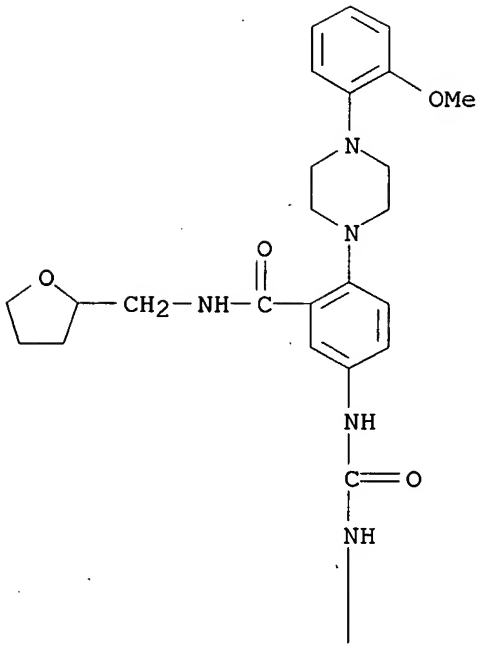
L4 ANSWER 5 OF 27 REGISTRY COPYRIGHT 2007 ACS on STN
RN 860078-73-5 REGISTRY
ED Entered STN: 12 Aug 2005
CN Pyrrolidine, 1-[5-[[[(3,5-dimethoxyphenyl)amino]carbonyl]amino]-2-[4-(2-methoxyphenyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)
MF C31 H37 N5 O5
SR Chemical Library



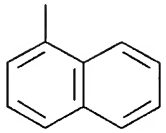
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 ANSWER 10 OF 27 REGISTRY COPYRIGHT 2007 ACS on STN
RN 857668-00-9 REGISTRY
ED Entered STN: 29 Jul 2005
CN Benzamide, 2-[4-(2-methoxyphenyl)-1-piperazinyl]-5-[[(1-naphthalenylamino)carbonyl]amino]-N-[(tetrahydro-2-furanyl)methyl]- (9CI) (CA INDEX NAME)
MF C34 H37 N5 O4
SR Chemical Library
Supplier: ComGenex International Inc.
LC STN Files: CHEMCATS

PAGE 1-A

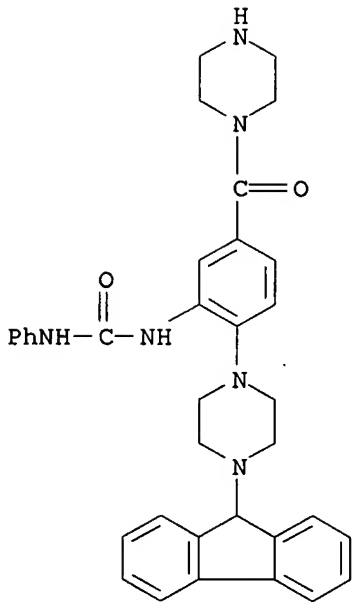


PAGE 2-A



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 ANSWER 27 OF 27 REGISTRY COPYRIGHT 2007 ACS on STN
RN 773882-48-7 REGISTRY
ED Entered STN: 02 Nov 2004
CN Piperazine, 1-[4-[4-(9H-fluoren-9-yl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)
MF C35 H36 N6 O2
SR CA
LC STN Files: CA, CAPLUS, USPATFULL



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> s l3 sss full
FULL SEARCH INITIATED 16:12:06 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 39190 TO ITERATE

100.0% PROCESSED 39190 ITERATIONS 422 ANSWERS
SEARCH TIME: 00.00.02

L5 422 SEA SSS FUL L3

=> save l5
ENTER NAME OR (END):ten815017/A
ANSWER SET L5 HAS BEEN SAVED AS 'TEN815017/A'

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	190.85	191.06

FILE 'CAPLUS' ENTERED AT 16:12:27 ON 28 APR 2007
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE COVERS 1907 - 28 Apr 2007 VOL 146 ISS 19
FILE LAST UPDATED: 27 Apr 2007 (20070427/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply.
They are available for your review at:

http://www.cas.org/infopolicy.html

=> s l5
L6 1 L5

=> d l6 1 bib abs fhitstr

L6 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2007 ACS on STN
AN 2004:857583 CAPLUS
DN 141:332220
TI A preparation of (piperazinylphenyl)urea derivatives as phospholipase C inhibitors, useful for the treatment of inflammatory disorders
IN Lagu, Bharat; Wachter, Michael; Rupert, Kenneth; Wachter, Michael
PA Janssen Pharmaceutica N.V., Belg.
SO PCT Int. Appl., 141 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004087685	A2	20041014	WO 2004-US9846	20040331
	WO 2004087685	A3	20041216		
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
	RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	US 2004235855	A1	20041125	US 2004-815017	20040331
PRAI	US 2003-458938P	P	20030331		

OS MARPAT 141:332220
GI

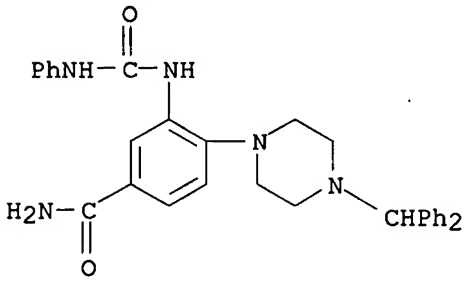
* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The invention relates to a preparation of (piperazinylphenyl)urea derivs. of

formula I [wherein: X is NH2, NH-alkyl, NHOH, NH-CN, or heterocyclic ring, etc.; Y is one or more (un)substituted alkyl; Z is (CH2)2-5; R1 is (un)substituted alkyl, cycloalkyl, or aryl, etc.; R2 is (un)substituted alkyl, C(O)alkyl, C(O)alkenyl, aryl, or cycloalkyl, etc.; R3 is O or S], useful as PLC-β2 inhibitors. For instance, (piperazinylphenyl)urea derivative II (IC50 = 1.2 μM) was prepared via addition of resin-bound (piperazinylphenyl)amine derivative III to Ph-N=C=O and subsequent resin cleavage (example 1).

IT 773882-10-3P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of (piperazinylphenyl)urea derivs. useful as PLC-β2 inhibitors)

RN 773882-10-3 CAPLUS
CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



=> file caold

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	5.74	196.80

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.78	-0.78

FILE 'CAOLD' ENTERED AT 16:13:01 ON 28 APR 2007
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE COVERS 1907-1966
FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, patent assignees, and patent information, e.g., patent numbers, are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

=> s l5
L7 0 L5

=> file chemcats

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.45	197.25

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.78

FILE 'CHEMCATS' ENTERED AT 16:13:11 ON 28 APR 2007
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE LAST UPDATED 21 APRIL 2007 (20070421/UP)

For details on recent updates in CHEMCATS, enter NEWS FILE at an arrow prompt. For the list of suppliers currently in the file, enter HELP SPA, HELP SPBC, HELP SPDH, HELP SPIN, HELP SPOP, and HELP SPQZ. For the list of current catalogs, enter HELP CTA, HELP CTBC, HELP CTDH, HELP CTIN, HELP CTOP, and HELP CTQZ.

This database is provided on an "as is" basis. Please consult the suppliers for current information regarding pricing, regional availability, available quantities, purities, etc. THERE ARE NO WARRANTIES OF ANY KIND, EITHER EXPRESSED OR IMPLIED. ACS is not liable for any loss of profit, goodwill or any other damages arising out of the use of this database.

CHEMCATS now contains more than 10 million records. See HELP CONTENT and NEWS FILE for details.

=> s l5
L8 287 L5

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.93	198.18

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.78

FILE 'REGISTRY' ENTERED AT 16:13:45 ON 28 APR 2007
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 27 APR 2007 HIGHEST RN 933069-51-3
DICTIONARY FILE UPDATES: 27 APR 2007 HIGHEST RN 933069-51-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

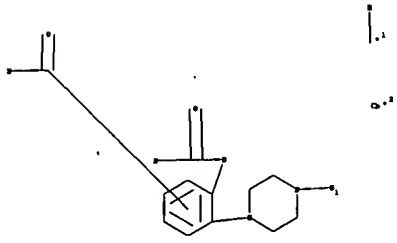
TSCA INFORMATION NOW CURRENT THROUGH December 2, 2006

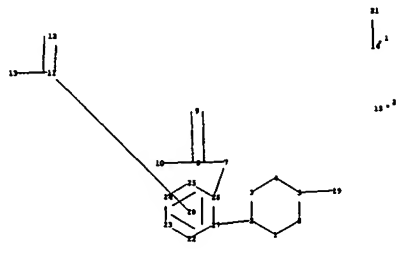
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

=>
Uploading C:\Documents and Settings\EBernhardt\My Documents\Stnexp\Queries\10815017-B.str





chain nodes :
7 8 9 10 11 12 14 15 19 21
ring nodes :
1 2 3 4 5 6 22 23 24 25 26 27
ring/chain nodes :

```
13
chain bonds :
2-27  5-19  7-8  7-26  8-9  8-10  11-12  11-13  14-21
ring bonds :
1-2  1-6  2-3  3-4  4-5  5-6  22-23  22-27  23-24  24-25  25-26  26-27
exact/norm bonds :
1-2  1-6  2-3  2-27  3-4  4-5  5-6  5-19  7-8  7-26  8-9  8-10  11-12  11-13
exact bonds :
14-21
normalized bonds :
22-23  22-27  23-24  24-25  25-26  26-27
isolated ring systems :
containing 1 :


G1:[*1],[*2]

Match level :
1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:CLASS  8:CLASS  9:CLASS  10:CLASS
 11:CLASS  12:CLASS  13:CLASS  14:CLASS  15:Atom  19:CLASS  21:CLASS  22:CLASS
23:Atom  24:Atom  25:Atom  26:CLASS  27:Atom  28:Atom



L9      STRUCTURE UPLOADED

=> s l9
SAMPLE SEARCH INITIATED 16:17:02 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -       36 TO ITERATE

100.0% PROCESSED       36 ITERATIONS                                 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:          360 TO     1080
PROJECTED ANSWERS:               1 TO       80

L10           1 SEA SSS SAM L9

=> d his

     (FILE 'HOME' ENTERED AT 16:07:18 ON 28 APR 2007)

     FILE 'REGISTRY' ENTERED AT 16:07:30 ON 28 APR 2007
L1              STRUCTURE UPLOADED
L2             34 S L1
L3              STRUCTURE UPLOADED
L4             27 S L3
L5            422 S L3 SSS FULL
                SAVE L5 TEN815017/A

     FILE 'CAPLUS' ENTERED AT 16:12:27 ON 28 APR 2007
L6             1 S L5

     FILE 'CAOLD' ENTERED AT 16:13:01 ON 28 APR 2007
L7             0 S L5
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     FILE 'CHEMCATS' ENTERED AT 16:13:11 ON 28 APR 2007
L8          287 S L5

     FILE 'REGISTRY' ENTERED AT 16:13:45 ON 28 APR 2007
L9              STRUCTURE UPLOADED
L10           1 S L9

=> s l9 sub=l5 full
FULL SUBSET SEARCH INITIATED 16:18:43 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED -       63 TO ITERATE

100.0% PROCESSED       63 ITERATIONS                                63 ANSWERS
SEARCH TIME: 00.00.01

L11           63 SEA SUB=L5 SSS FUL L9

=> file caplus
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                         ENTRY    SESSION
FULL ESTIMATED COST                                      44.25     242.43

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)          SINCE FILE      TOTAL
                                                         ENTRY    SESSION
CA SUBSCRIBER PRICE                                       0.00      -0.78

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FILE LAST UPDATED: 27 Apr 2007  (20070427/ED)

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=> s l11
L12           1 L11

=> d l12 bib abs hitstr

L12  ANSWER 1 OF 1  CAPLUS  COPYRIGHT 2007 ACS on STN
AN   2004:857583  CAPLUS
DN   141:332220
```

TI A preparation of (piperazinylphenyl)urea derivatives as phospholipase C inhibitors, useful for the treatment of inflammatory disorders
IN Lagu, Bharat; Wachter, Michael; Rupert, Kenneth; Wachter, Michael
PA Janssen Pharmaceutica N.V., Belg.
SO PCT Int. Appl., 141 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2004087685	A2	20041014	WO 2004-US9846	20040331
WO 2004087685	A3	20041216		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
US 2004235855	A1	20041125	US 2004-815017	20040331
PRAI US 2003-458938P	P	20030331		

OS MARPAT 141:332220
GI

Apps

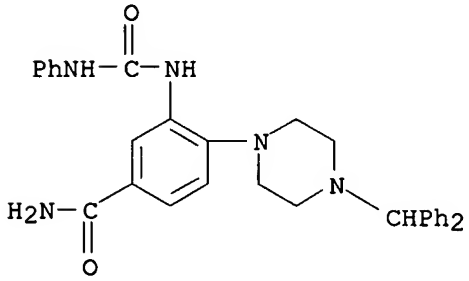
* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The invention relates to a preparation of (piperazinylphenyl)urea derivs. of formula I [wherein: X is NH2, NH-alkyl, NHOH, NH-CN, or heterocyclic ring, etc.; Y is one or more (un)substituted alkyl; Z is (CH2)2-5; R1 is (un)substituted alkyl, cycloalkyl, or aryl, etc.; R2 is (un)substituted alkyl, C(O)alkyl, C(O)alkenyl, aryl, or cycloalkyl, etc.; R3 is O or S], useful as PLC-β2 inhibitors. For instance, (piperazinylphenyl)urea derivative II (IC50 = 1.2 μM) was prepared via addition of resin-bound (piperazinylphenyl)amine derivative III to Ph-N=C=O and subsequent resin cleavage (example 1).

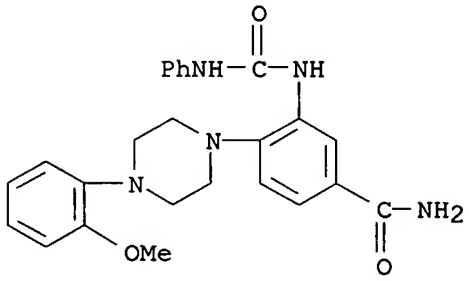
IT 773882-10-3P 773882-11-4P 773882-12-5P
773882-13-6P 773882-14-7P 773882-15-8P
773882-16-9P 773882-17-0P 773882-18-1P
773882-19-2P 773882-20-5P 773882-21-6P
773882-22-7P 773882-23-8P 773882-24-9P
773882-25-0P 773882-26-1P 773882-27-2P
773882-28-3P 773882-29-4P 773882-32-9P
773882-33-0P 773882-34-1P 773882-35-2P
773882-36-3P 773882-37-4P 773882-38-5P
773882-39-6P 773882-40-9P 773882-41-0P
773882-42-1P 773882-43-2P 773882-44-3P
773882-45-4P 773882-46-5P 773882-47-6P
773882-48-7P 773882-49-8P 773882-50-1P
773882-52-3P 773882-54-5P 773882-56-7P
773882-58-9P 773882-59-0P 773882-61-4P

773882-63-6P 773882-65-8P 773882-66-9P
773882-67-0P 773882-68-1P 773882-69-2P
773882-70-5P 773882-71-6P 773882-72-7P
773882-73-8P 773882-74-9P 773882-76-1P
773882-78-3P 773882-79-4P 773882-85-2P
773882-86-3P 773882-87-4P 773882-88-5P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of (piperazinylphenyl)urea derivs. useful as PLC-β2 inhibitors)

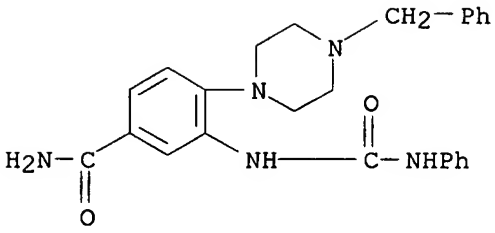
RN 773882-10-3 CAPLUS
CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]-. (9CI) (CA INDEX NAME)



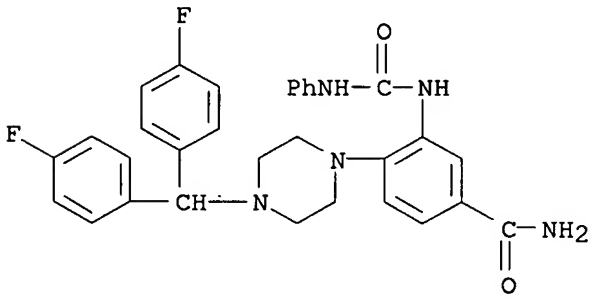
RN 773882-11-4 CAPLUS
CN Benzamide, 4-[4-(2-methoxyphenyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



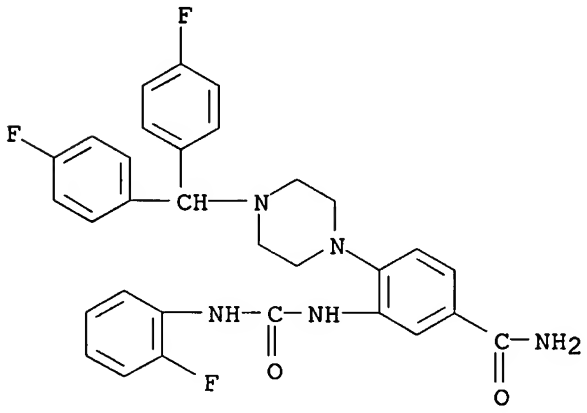
RN 773882-12-5 CAPLUS
CN Benzamide, 3-[[(phenylamino)carbonyl]amino]-4-[4-(phenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



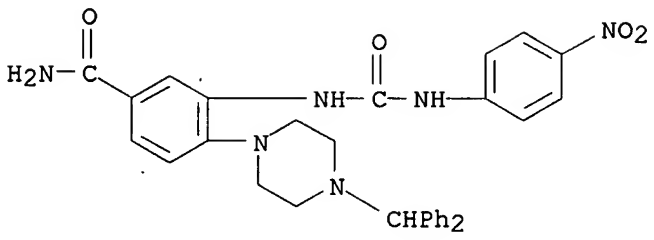
RN 773882-13-6 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



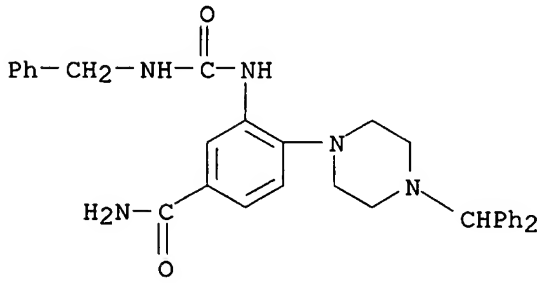
RN 773882-14-7 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(2-fluorophenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



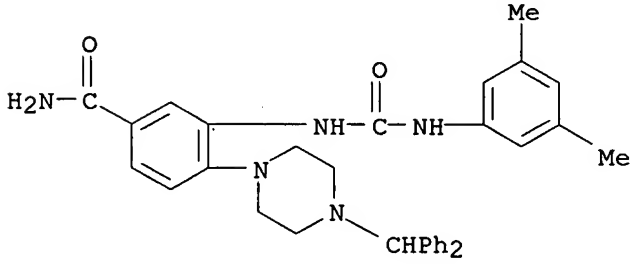
RN 773882-15-8 CAPLUS
CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(4-nitrophenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



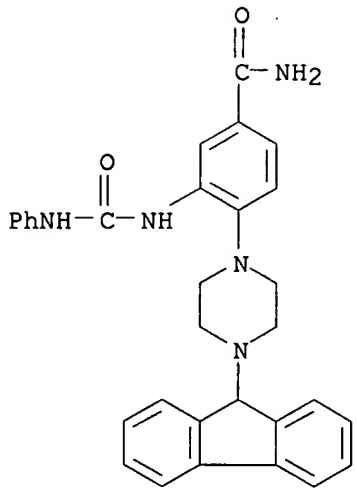
RN 773882-16-9 CAPLUS
CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(phenylmethyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



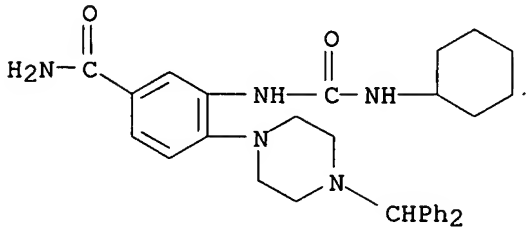
RN 773882-17-0 CAPLUS
CN Benzamide, 3-[[[(3,5-dimethylphenyl)amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



RN 773882-18-1 CAPLUS
CN Benzamide, 4-[4-(9H-fluoren-9-yl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)

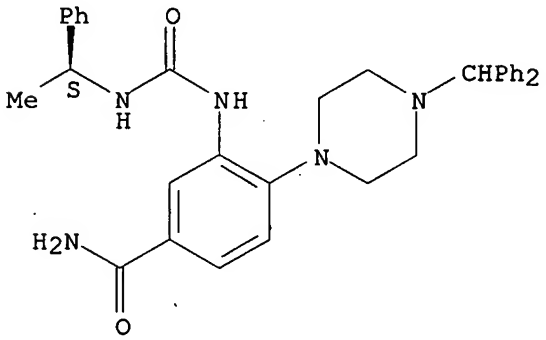


RN 773882-19-2 CAPLUS
CN Benzamide, 3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)

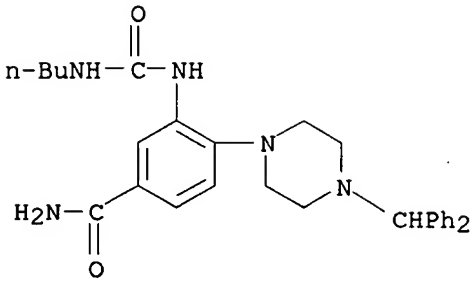


RN 773882-20-5 CAPLUS
CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[[(1S)-1-phenylethyl]amino]carbonyl]amino]- (9CI) (CA INDEX NAME)

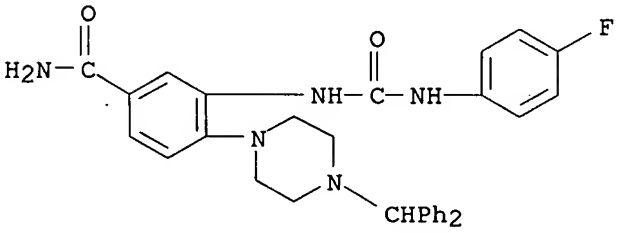
Absolute stereochemistry.



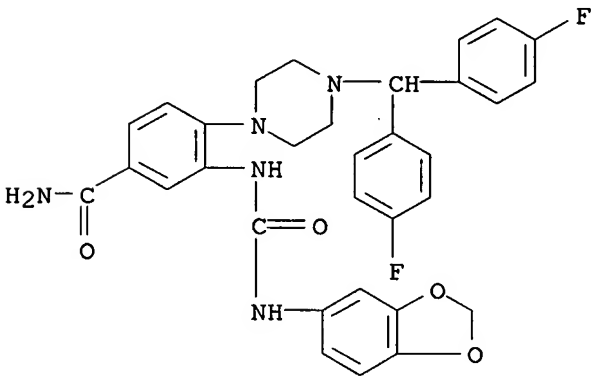
RN 773882-21-6 CAPLUS
CN Benzamide, 3-[[(butylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



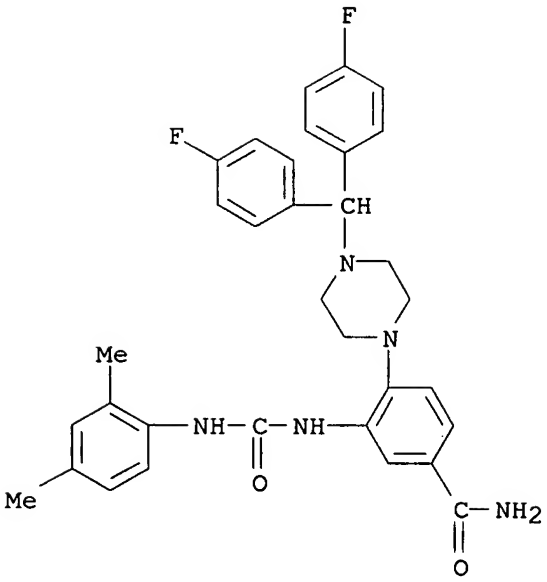
RN 773882-22-7 CAPLUS
CN Benzamide, 4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(4-fluorophenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



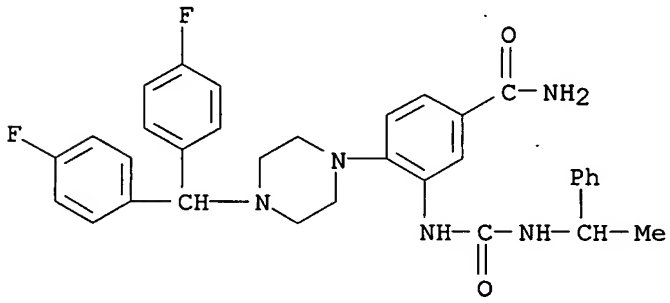
RN 773882-23-8 CAPLUS
CN Benzamide, 3-[[(1,3-benzodioxol-5-ylamino)carbonyl]amino]-4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)



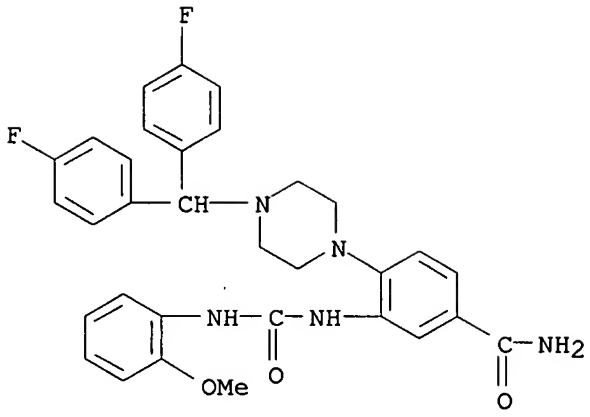
RN 773882-24-9 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(2,4-dimethylphenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



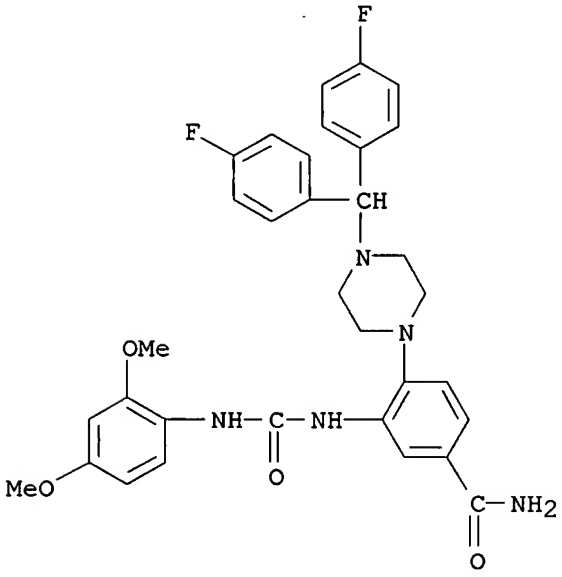
RN 773882-25-0 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(1-phenylethyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



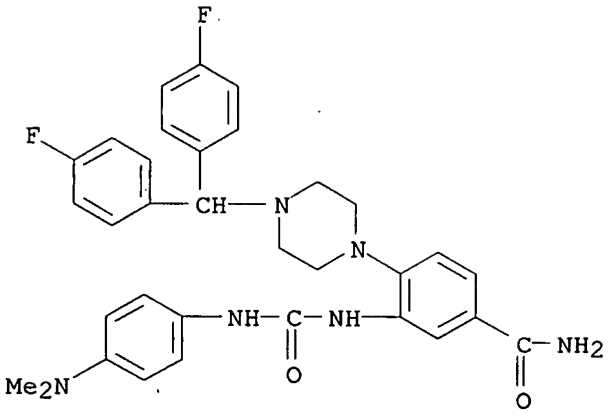
RN 773882-26-1 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(2-methoxyphenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



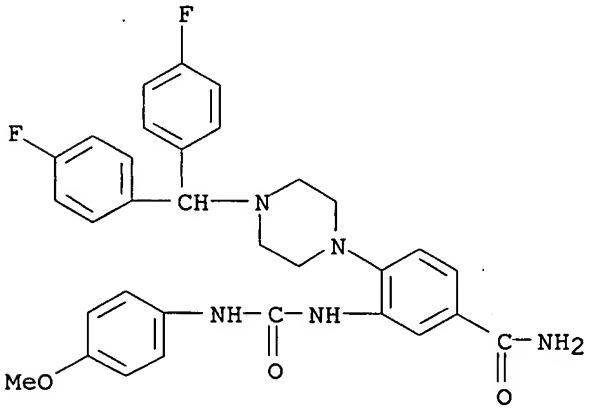
RN 773882-27-2 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(2,4-dimethoxyphenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



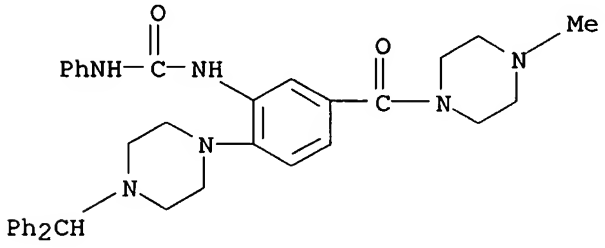
RN 773882-28-3 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[[4-(dimethylamino)phenyl]amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



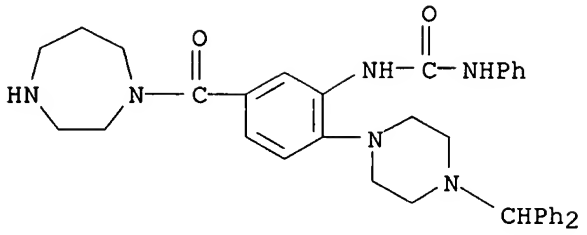
RN 773882-29-4 CAPLUS
CN Benzamide, 4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[[(4-methoxyphenyl)amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



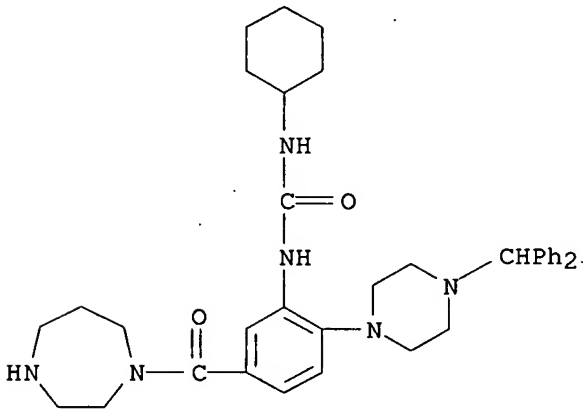
RN 773882-32-9 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]-4-methyl- (9CI) (CA INDEX NAME)



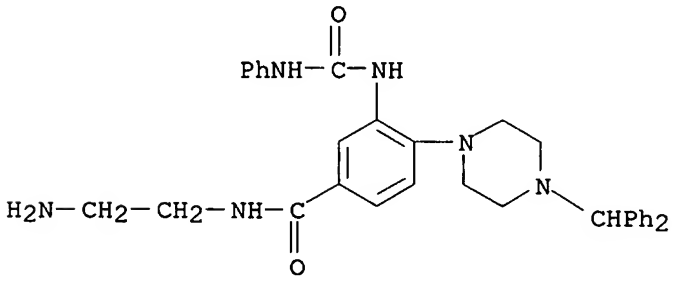
RN 773882-33-0 CAPLUS
CN 1H-1,4-Diazepine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]hexahydro- (9CI) (CA INDEX NAME)



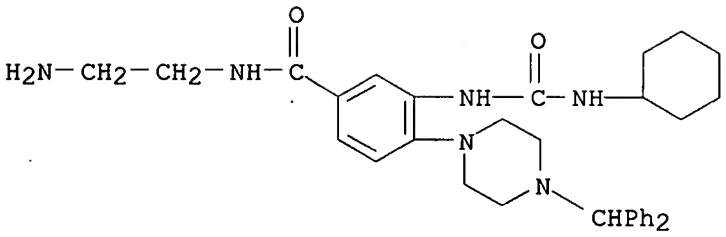
RN 773882-34-1 CAPLUS
CN 1H-1,4-Diazepine, 1-[3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]hexahydro- (9CI) (CA INDEX NAME)



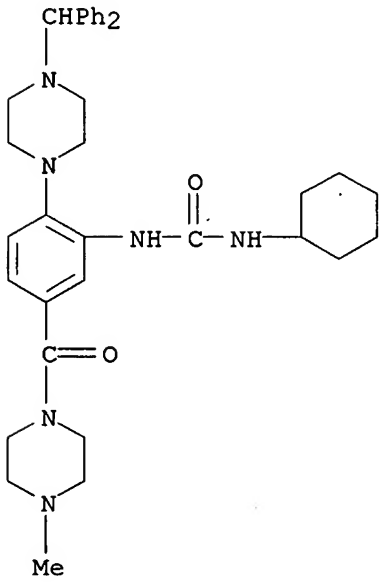
RN 773882-35-2 CAPLUS
CN Benzamide, N-(2-aminoethyl)-4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



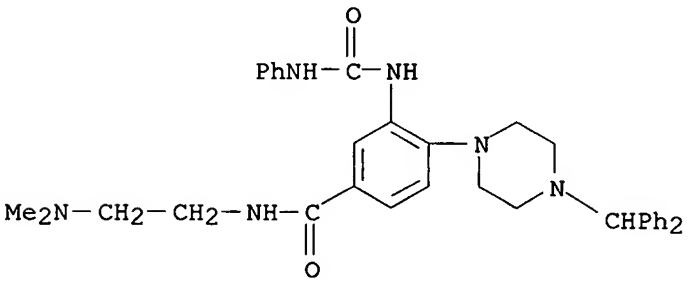
RN 773882-36-3 CAPLUS
CN Benzamide, N-(2-aminoethyl)-3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



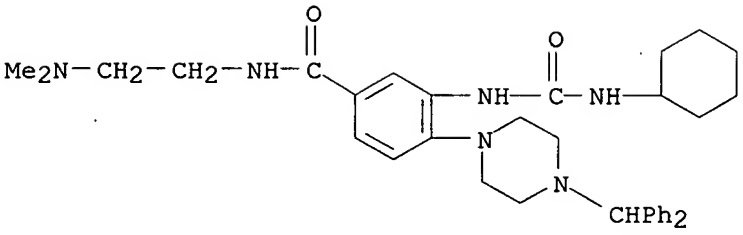
RN 773882-37-4 CAPLUS
CN Piperazine, 1-[3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]-4-methyl- (9CI) (CA INDEX NAME)



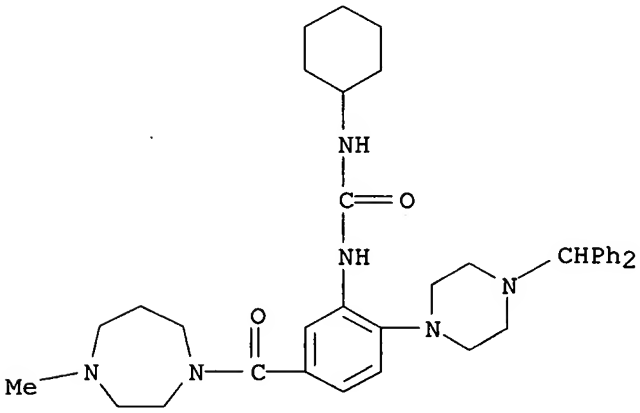
RN 773882-38-5 CAPLUS
CN Benzamide, N-[2-(dimethylamino)ethyl]-4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



RN 773882-39-6 CAPLUS
CN Benzamide, 3-[[(cyclohexylamino)carbonyl]amino]-N-[2-(dimethylamino)ethyl]-4-[4-(diphenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)

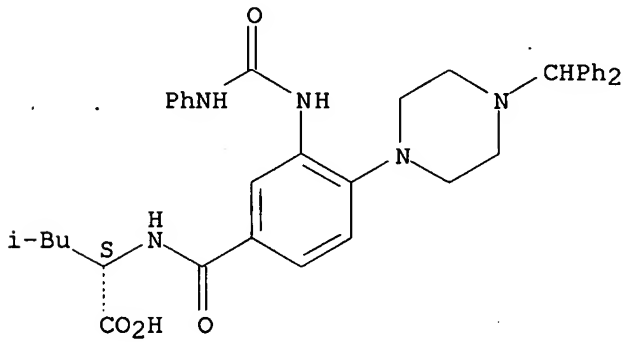


RN 773882-40-9 CAPLUS
CN 1H-1,4-Diazepine, 1-[3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]hexahydro-4-methyl- (9CI) (CA INDEX NAME)

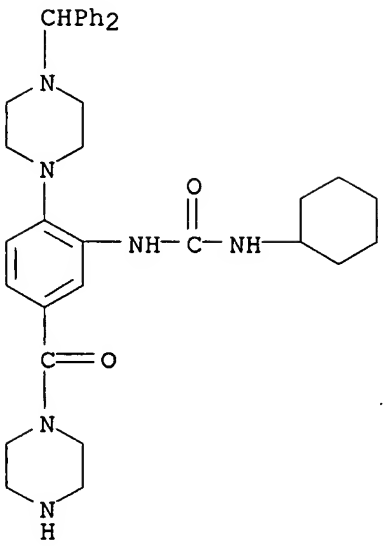


RN 773882-41-0 CAPLUS
CN L-Leucine, N-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)

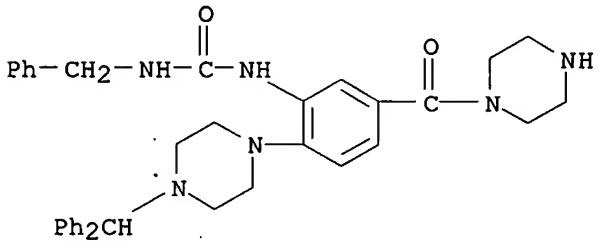
Absolute stereochemistry.



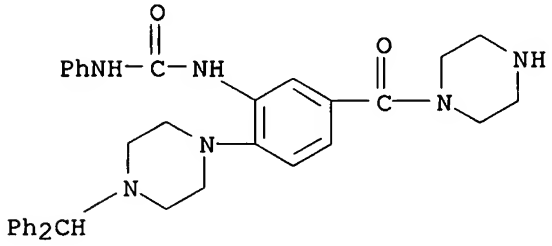
RN 773882-42-1 CAPLUS
CN Piperazine, 1-[3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



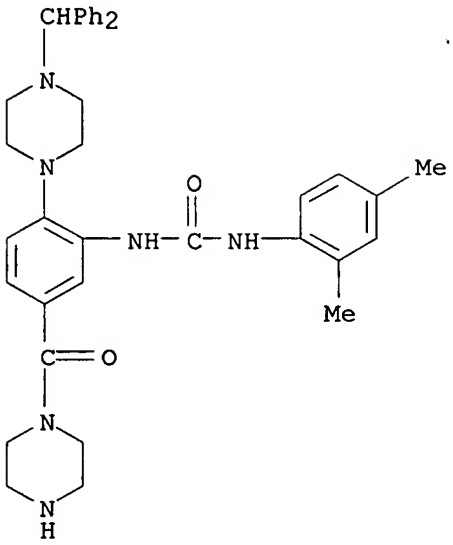
RN 773882-43-2 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(phenylmethyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



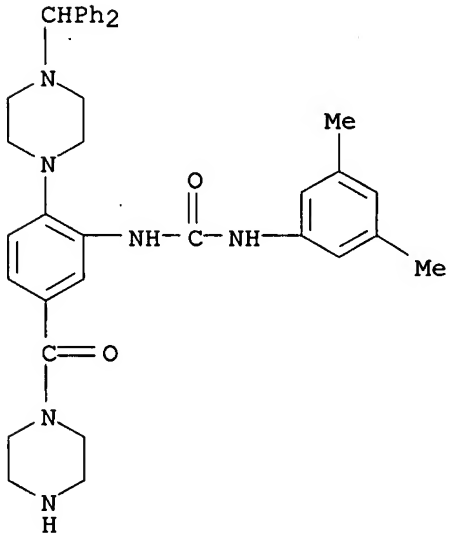
RN 773882-44-3 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



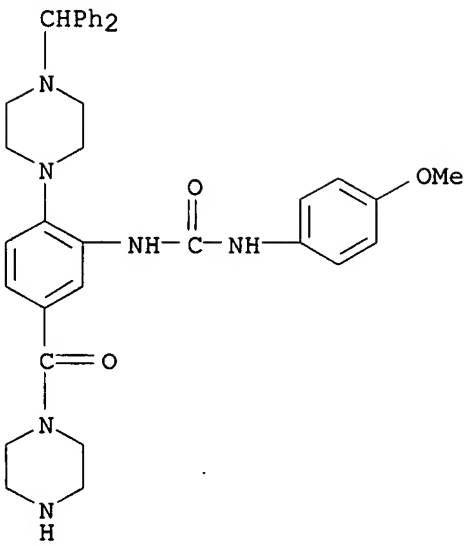
RN 773882-45-4 CAPLUS
CN Piperazine, 1-[3-[[[(2,4-dimethylphenyl)amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



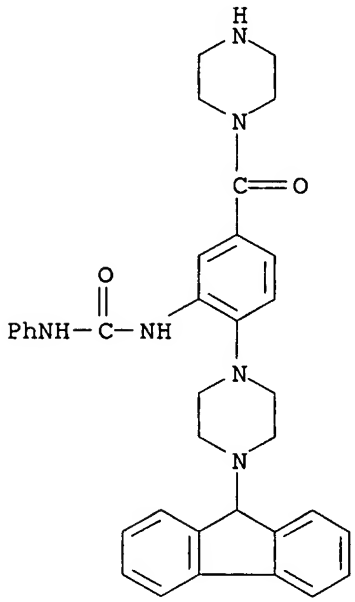
RN 773882-46-5 CAPLUS
CN Piperazine, 1-[3-[[[(3,5-dimethylphenyl)amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



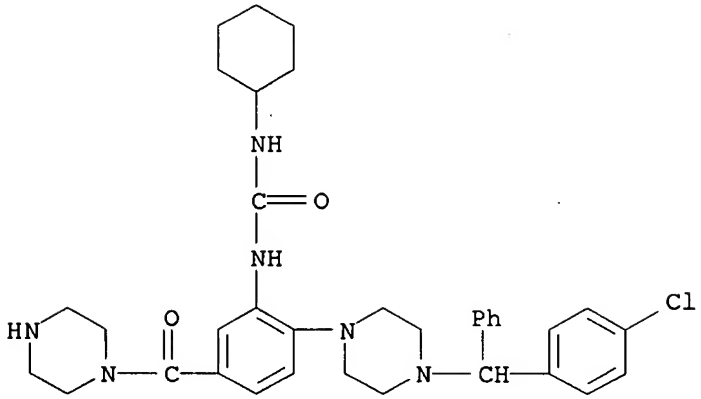
RN 773882-47-6 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(4-methoxyphenyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



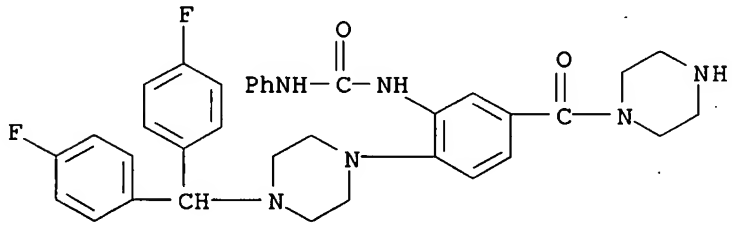
RN 773882-48-7 CAPLUS
CN Piperazine, 1-[4-[4-(9H-fluoren-9-yl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



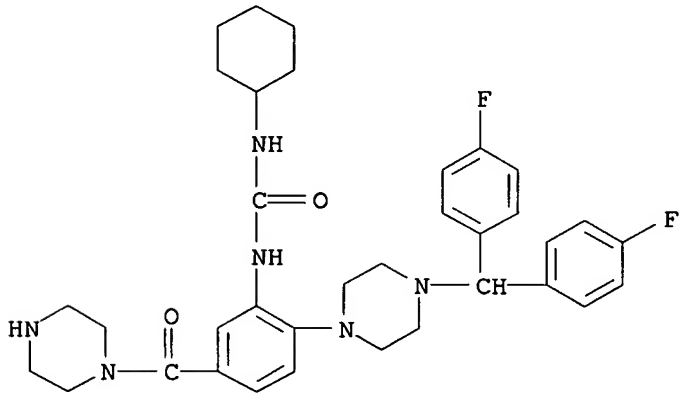
RN 773882-49-8 CAPLUS
CN Piperazine, 1-[4-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]-3-[[(cyclohexylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



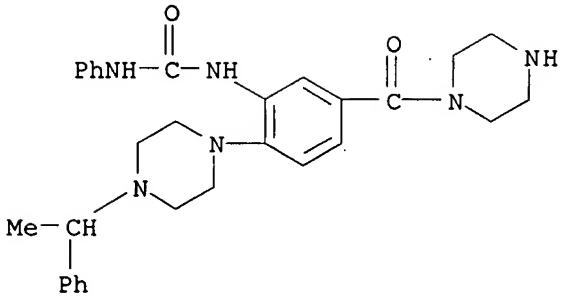
RN 773882-50-1 CAPLUS
CN Piperazine, 1-[4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



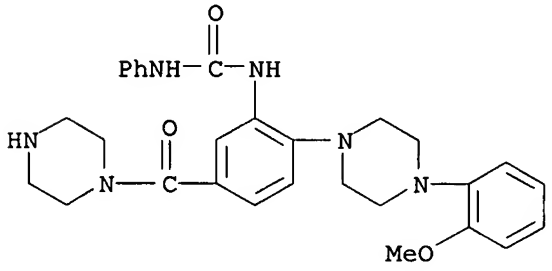
RN 773882-52-3 CAPLUS
CN Piperazine, 1-[4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[(cyclohexylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



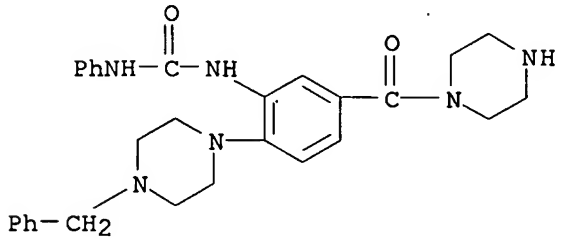
RN 773882-54-5 CAPLUS
CN Piperazine, 1-[3-[[(phenylamino)carbonyl]amino]-4-[4-(1-phenylethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



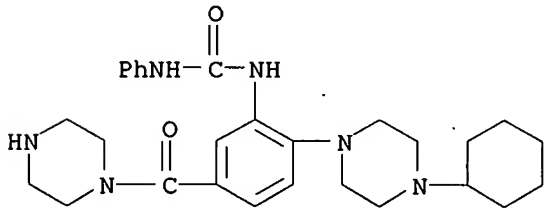
RN 773882-56-7 CAPLUS
CN Piperazine, 1-[4-[4-(2-methoxyphenyl)-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



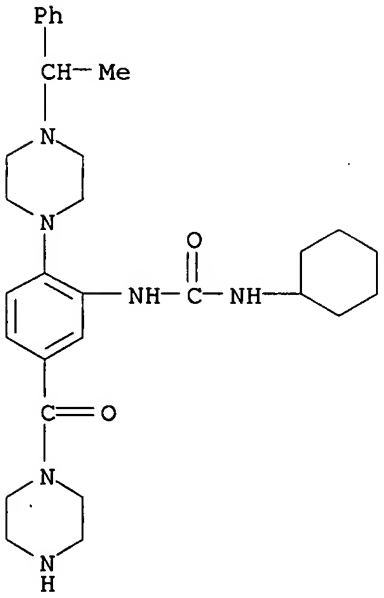
RN 773882-58-9 CAPLUS
CN Piperazine, 1-[3-[[(phenylamino)carbonyl]amino]-4-[4-(phenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



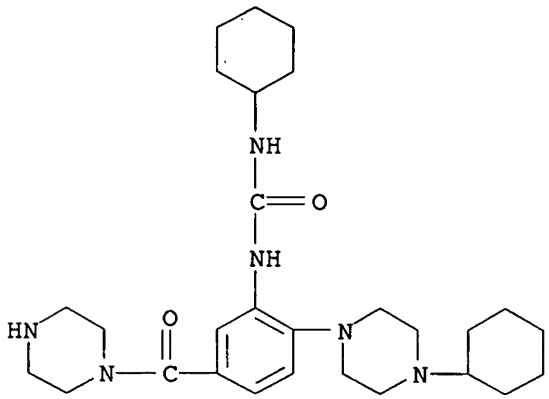
RN 773882-59-0 CAPLUS
CN Piperazine, 1-[4-(4-cyclohexyl-1-piperazinyl)-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



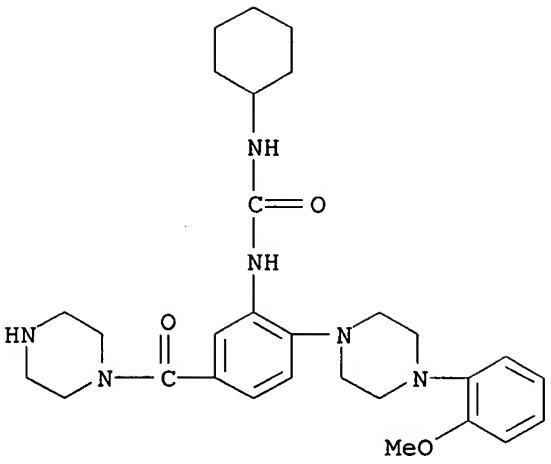
RN 773882-61-4 CAPLUS
CN Piperazine, 1-[3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(1-phenylethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



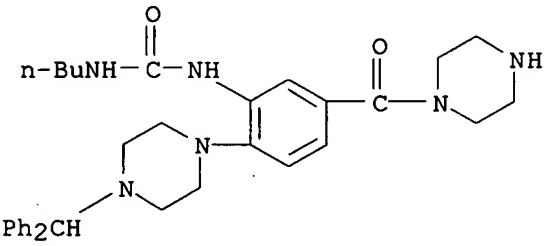
RN 773882-63-6 CAPLUS
CN Piperazine, 1-[3-[[(cyclohexylamino)carbonyl]amino]-4-(4-cyclohexyl-1-piperazinyl)benzoyl]- (9CI) (CA INDEX NAME)



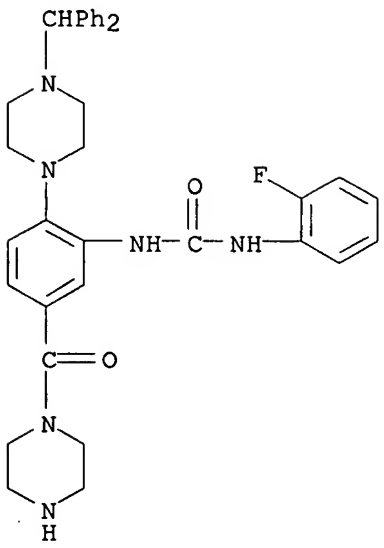
RN 773882-65-8 CAPLUS
CN Piperazine, 1-[3-[[(cyclohexylamino)carbonyl]amino]-4-[4-(2-methoxyphenyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



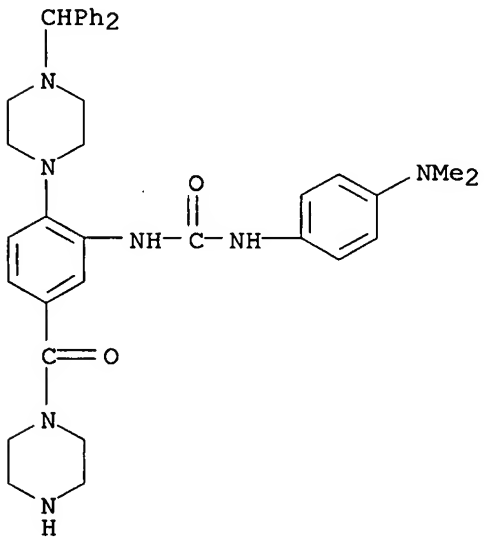
RN 773882-66-9 CAPLUS
CN Piperazine, 1-[3-[[(butylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



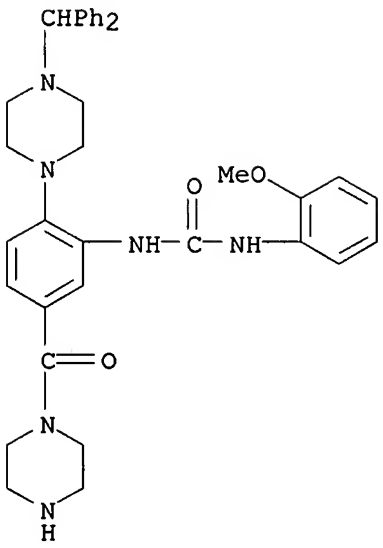
RN 773882-67-0 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(2-fluorophenyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



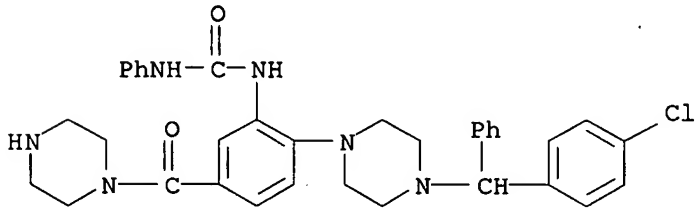
RN 773882-68-1 CAPLUS
CN Piperazine, 1-[3-[[[[4-(dimethylamino)phenyl]amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



RN 773882-69-2 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(2-methoxyphenyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)

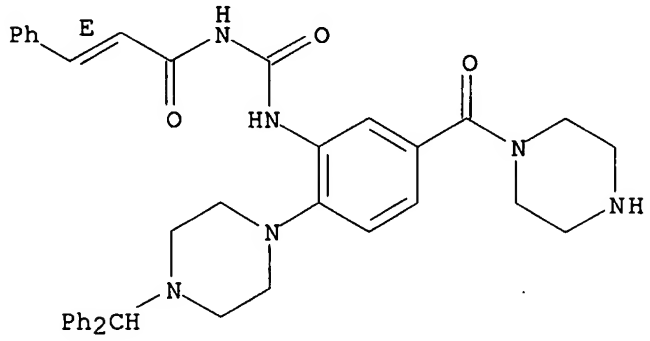


RN 773882-70-5 CAPLUS
CN Piperazine, 1-[4-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



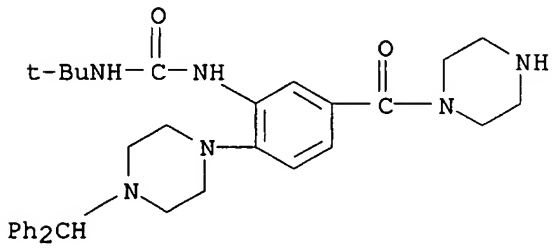
RN 773882-71-6 CAPLUS
CN 2-Propenamide, N-[[[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]amino]carbonyl]-3-phenyl-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

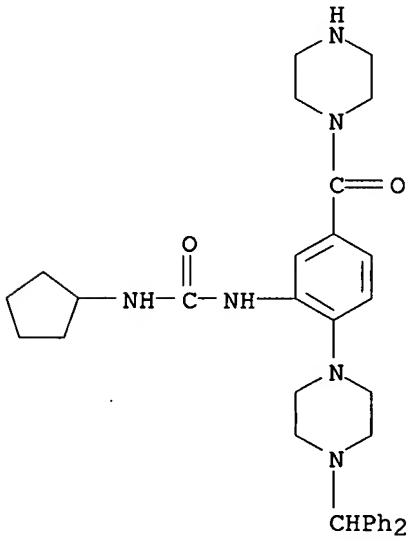


RN 773882-72-7 CAPLUS

CN Piperazine, 1-[3-[[[(1,1-dimethylethyl)amino]carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)

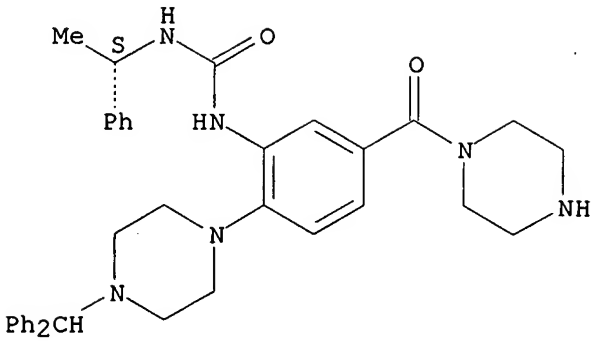


RN 773882-73-8 CAPLUS
CN Piperazine, 1-[3-[[(cyclopentylamino)carbonyl]amino]-4-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)

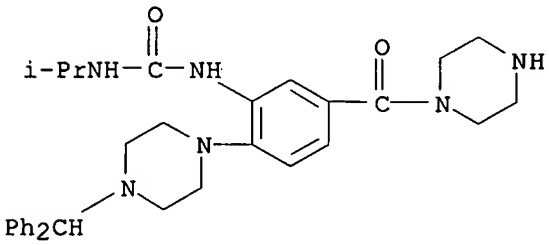


RN 773882-74-9 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[[(1S)-1-phenylethyl]amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)

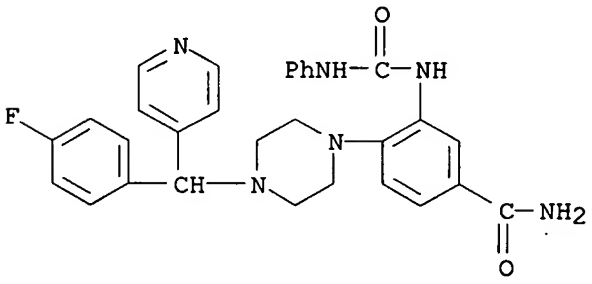
Absolute stereochemistry.



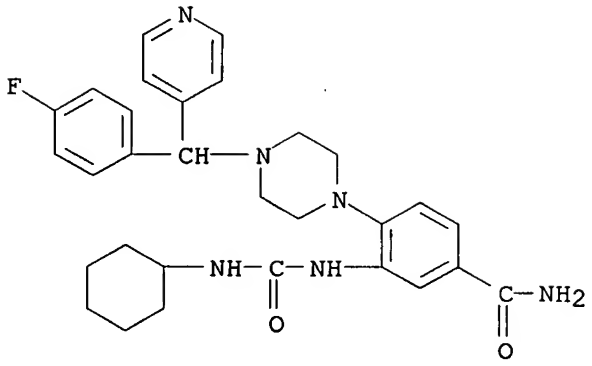
RN 773882-76-1 CAPLUS
CN Piperazine, 1-[4-[4-(diphenylmethyl)-1-piperazinyl]-3-[[[(1-methylethyl)amino]carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



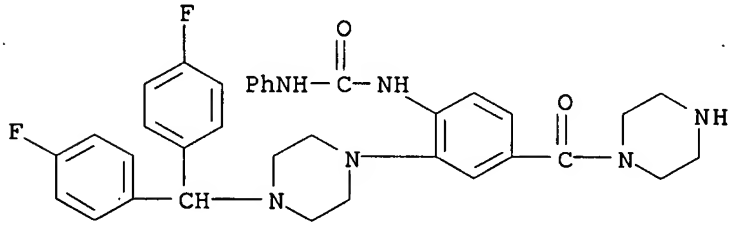
RN 773882-78-3 CAPLUS
CN Benzamide, 4-[4-[(4-fluorophenyl)-4-pyridinylmethyl]-1-piperazinyl]-3-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



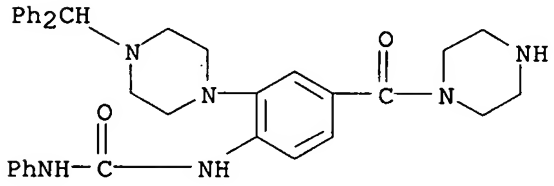
RN 773882-79-4 CAPLUS
CN Benzamide, 3-[[(cyclohexylamino)carbonyl]amino]-4-[4-[(4-fluorophenyl)-4-pyridinylmethyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)



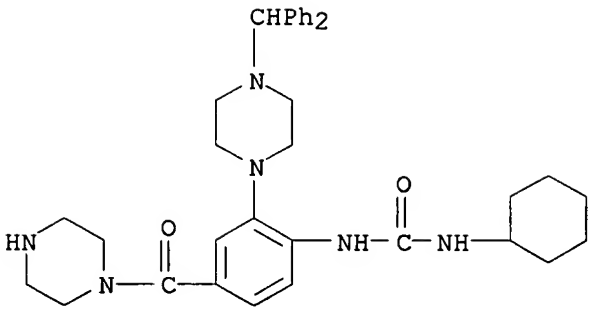
RN 773882-85-2 CAPLUS
CN Piperazine, 1-[3-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-4-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



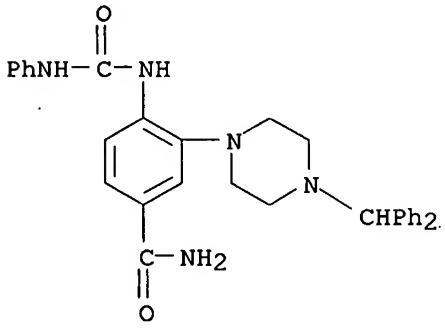
RN 773882-86-3 CAPLUS
CN Piperazine, 1-[3-[4-(diphenylmethyl)-1-piperazinyl]-4-[[(phenylamino)carbonyl]amino]benzoyl]- (9CI) (CA INDEX NAME)



RN 773882-87-4 CAPLUS
CN Piperazine, 1-[4-[[(cyclohexylamino)carbonyl]amino]-3-[4-(diphenylmethyl)-1-piperazinyl]benzoyl]- (9CI) (CA INDEX NAME)



RN 773882-88-5 CAPLUS
CN Benzamide, 3-[4-(diphenylmethyl)-1-piperazinyl]-4-[[(phenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



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